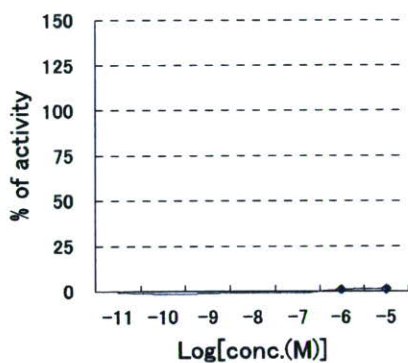


sample No. AT113  
 chemical name Uracil, 6-propyl-2-thio-  
 CAS. 51-52-5

**AR agonist assay**

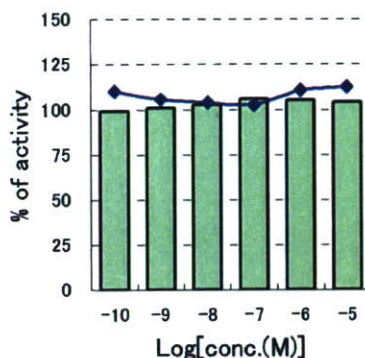
LOG[conc.(M)]	luc
-5	1.1
-6	1.0
-7	-0.9
-8	-0.7
-9	-1.1
-10	-1.6
-11	-1.0



**AR antagonist assay**

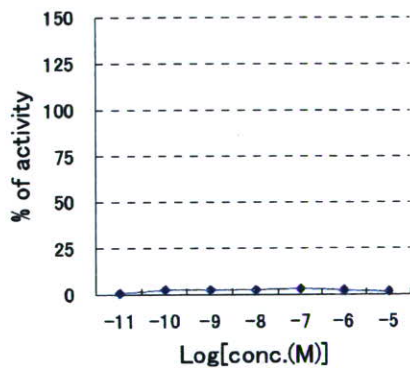
LOG[conc.(M)]	luc	ren
-5	113	104
-6	111	105
-7	103	106
-8	104	102
-9	106	101
-10	110	99

Cell viability



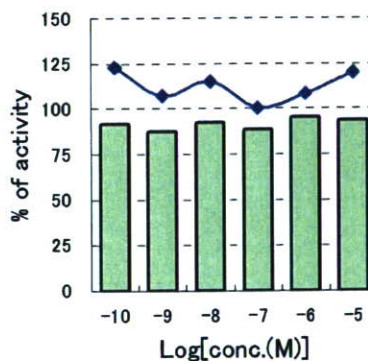
**TRβ-RXR agonist assay**

LOG[conc.(M)]	luc
-5	1.8
-6	2.5
-7	3.4
-8	2.8
-9	2.7
-10	2.7
-11	1.0



**TRβ-RXR antagonist assay**

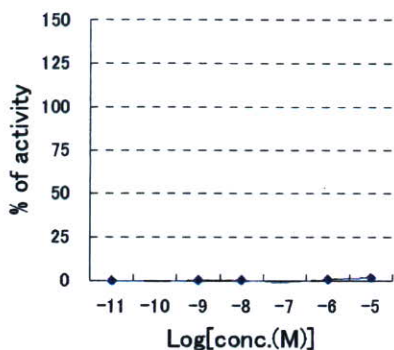
LOG[conc.(M)]	luc	ren
-5	120	93
-6	109	95
-7	101	89
-8	115	92
-9	107	88
-10	123	92



sample No. AT114  
 chemical name 3-Amino-s-triazole  
 CAS. 61-82-5

**AR agonist assay**

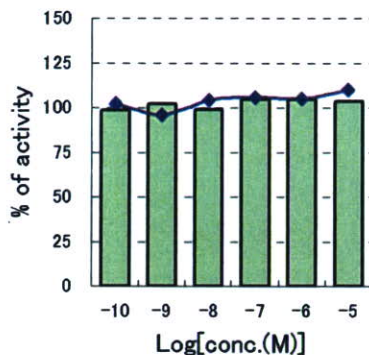
LOG[conc.(M)]	luc
-5	1.5
-6	0.6
-7	-1.1
-8	0.2
-9	0.2
-10	-0.6
-11	-0.1



**AR antagonist assay**

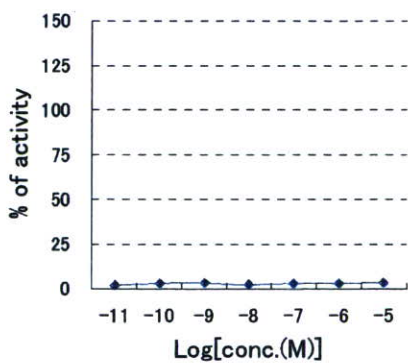
LOG[conc.(M)]	luc	ren
-5	110	104
-6	105	104
-7	106	105
-8	104	99
-9	96	102
-10	102	99

Cell viability



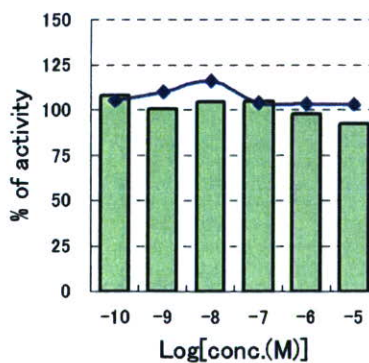
**TRβ-RXR agonist assay**

LOG[conc.(M)]	luc
-5	3.6
-6	3.1
-7	3.1
-8	2.6
-9	3.5
-10	3.1
-11	2.4



**TRβ-RXR antagonist assay**

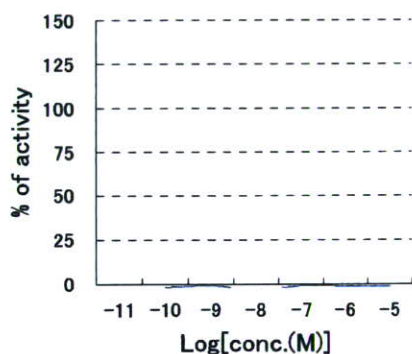
LOG[conc.(M)]	luc	ren
-5	103	92
-6	103	98
-7	104	105
-8	116	104
-9	110	101
-10	106	108



sample No. AT115  
 chemical name Pentadecanoic acid, 15-hydroxy-, .xi.-lactone  
 CAS. 106-02-5

**AR agonist assay**

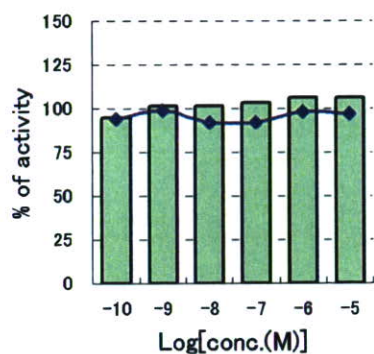
LOG[conc.(M)]	luc
-5	-1.2
-6	-1.2
-7	-0.9
-8	-3.3
-9	-1.1
-10	-1.8
-11	-2.4



**AR antagonist assay**

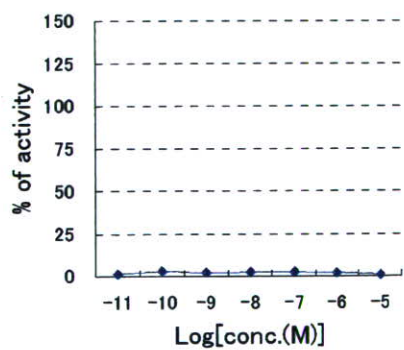
LOG[conc.(M)]	luc	ren
-5	97	106
-6	98	106
-7	92	103
-8	92	101
-9	99	101
-10	94	95

Cell viability



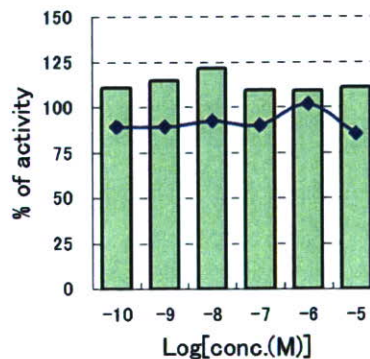
**TRβ-RXR agonist assay**

LOG[conc.(M)]	luc
-5	0.9
-6	2.0
-7	2.7
-8	2.7
-9	2.2
-10	3.1
-11	1.2



**TRβ-RXR antagonist assay**

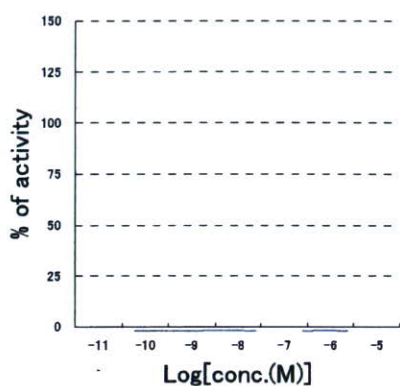
LOG[conc.(M)]	luc	ren
-5	86	111
-6	102	109
-7	91	110
-8	93	122
-9	89	115
-10	89	111



sample No. AT116  
 chemical name Morpholine, 4,4'-dithiodi-  
 CAS. 103-34-4

**AR agonist assay**

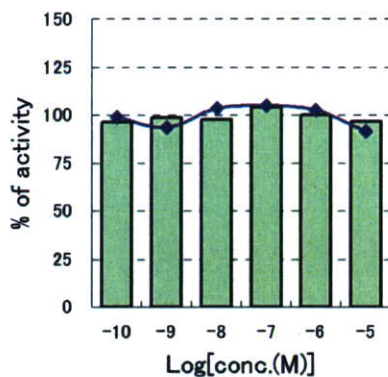
LOG[conc.(M)]	luc
-5	-2.7
-6	-1.5
-7	-2.4
-8	-1.4
-9	-1.7
-10	-1.8
-11	-2.2



**AR antagonist assay**

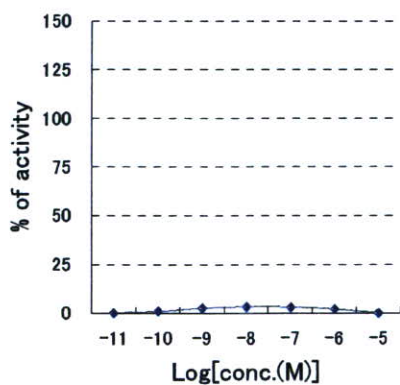
LOG[conc.(M)]	luc	ren
-5	92	97
-6	103	100
-7	105	104
-8	104	98
-9	94	98
-10	99	96

Cell viability



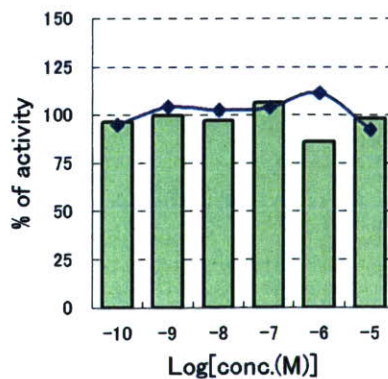
**TRβ-RXR agonist assay**

LOG[conc.(M)]	luc
-5	0.3
-6	2.3
-7	3.4
-8	3.5
-9	2.8
-10	1.2
-11	0.3



**TRβ-RXR antagonist assay**

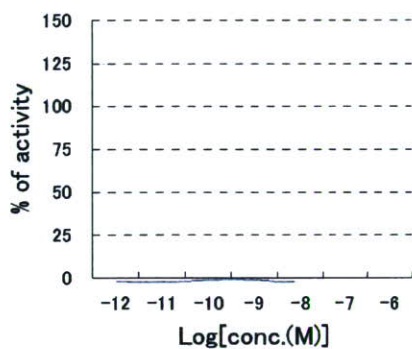
LOG[conc.(M)]	luc	ren
-5	92	98
-6	111	86
-7	104	106
-8	103	97
-9	104	100
-10	95	96



sample No. AT117  
 chemical name 7H-Furo(3,2-g)(1)benzopyran-7-one, 9-methoxy-  
 CAS. 298-81-7

**AR agonist assay**

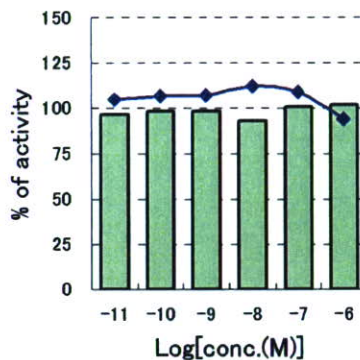
LOG[conc.(M)]	luc
-6	-1.9
-7	-1.9
-8	-2.6
-9	-1.1
-10	-0.9
-11	-2.2
-12	-1.6



**AR antagonist assay**

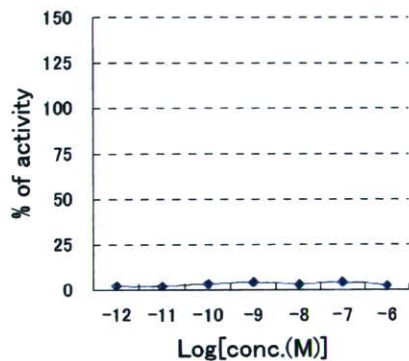
LOG[conc.(M)]	luc	ren
-6	94	102
-7	109	101
-8	112	93
-9	107	99
-10	107	98
-11	105	97

Cell viability



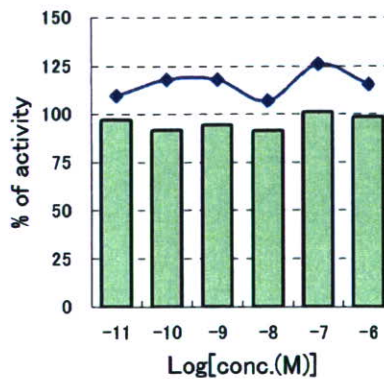
**TRβ-RXR agonist assay**

LOG[conc.(M)]	luc
-6	2.5
-7	4.5
-8	3.2
-9	4.6
-10	3.6
-11	2.2
-12	2.6



**TRβ-RXR antagonist assay**

LOG[conc.(M)]	luc	ren
-6	115	98
-7	126	101
-8	107	91
-9	118	94
-10	118	92
-11	110	97

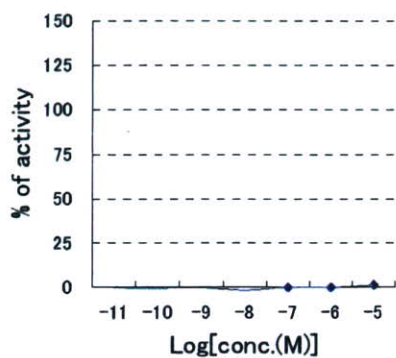




sample No. AT118  
 chemical name Fenvalerate  
 CAS. 51630-58-1

**AR agonist assay**

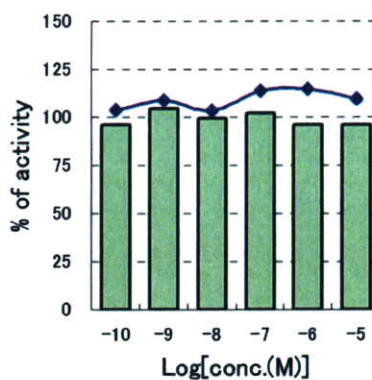
LOG[conc.(M)]	luc
-5	1.3
-6	-0.1
-7	-0.1
-8	-1.6
-9	-0.3
-10	-0.7
-11	-0.2



**AR antagonist assay**

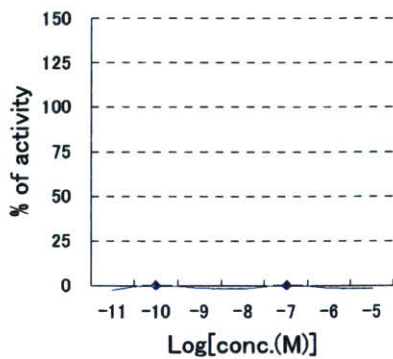
LOG[conc.(M)]	luc	ren
-5	110	96
-6	115	96
-7	114	102
-8	103	99
-9	109	104
-10	104	96

Cell viability



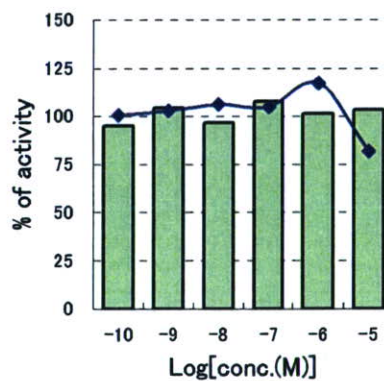
**TRβ-RXR agonist assay**

LOG[conc.(M)]	luc
-5	-1.3
-6	-1.0
-7	0.6
-8	-1.3
-9	-1.0
-10	0.6
-11	-1.8



**TRβ-RXR antagonist assay**

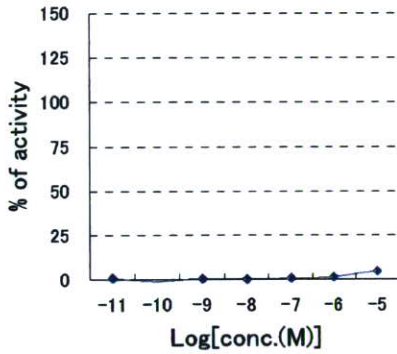
LOG[conc.(M)]	luc	ren
-5	82	103
-6	117	101
-7	105	108
-8	106	96
-9	103	104
-10	101	95



sample No. AT119  
 chemical name Anthraquinone, 1,4-dihydroxy-  
 CAS. 81-64-1

**AR agonist assay**

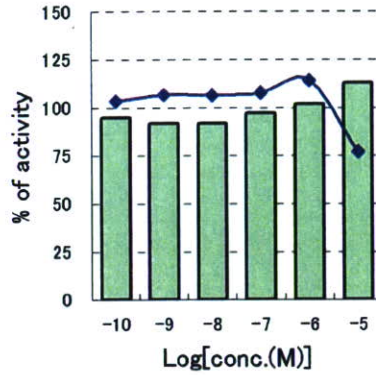
LOG[conc.(M)]	luc
-5	4.8
-6	1.7
-7	0.9
-8	0.5
-9	0.8
-10	-1.1
-11	1.0



**AR antagonist assay**

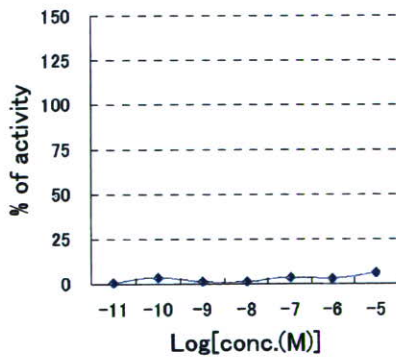
LOG[conc.(M)]	luc	ren
-5	77	113
-6	114	102
-7	108	97
-8	107	92
-9	107	92
-10	104	95

Cell viability



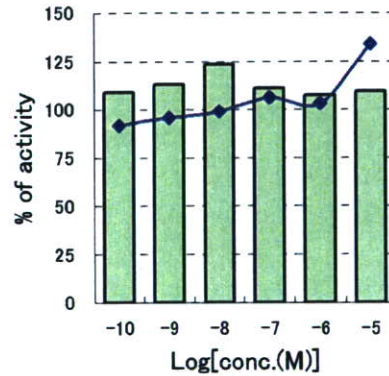
**TRβ-RXR agonist assay**

LOG[conc.(M)]	luc
-5	6.6
-6	3.4
-7	3.9
-8	1.6
-9	1.5
-10	4.0
-11	1.1



**TRβ-RXR antagonist assay**

LOG[conc.(M)]	luc	ren
-5	134	109
-6	103	107
-7	106	111
-8	99	123
-9	96	113
-10	92	109

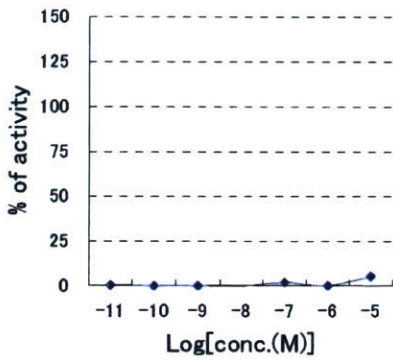


sample No.  
chemical name  
CAS.

AT120  
Phenol, 4,4'-(1-(4-(1-(4-hydroxyphenyl)-1-methylethyl)phenyl)ethylidene)bis-  
110726-28-8

**AR agonist assay**

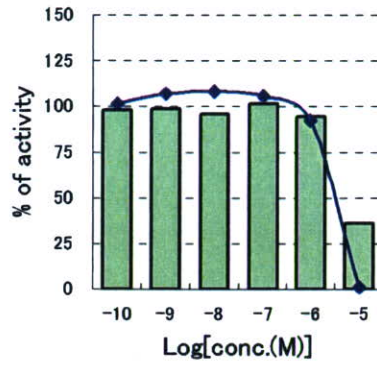
LOG[conc.(M)]	luc
-5	5.4
-6	0.1
-7	2.0
-8	-0.4
-9	0.0
-10	0.0
-11	0.2



**AR antagonist assay**

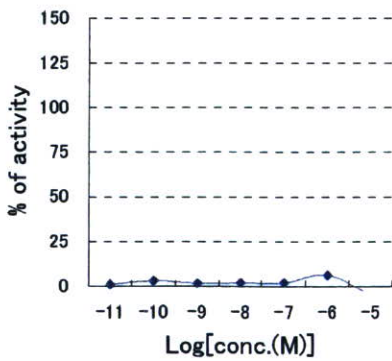
LOG[conc.(M)]	luc	ren
-5	1	36
-6	92	95
-7	106	102
-8	108	96
-9	107	99
-10	102	98

Cell viability



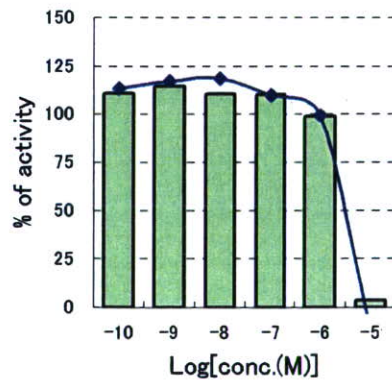
**TRβ-RXR agonist assay**

LOG[conc.(M)]	luc
-5	-4.3
-6	6.6
-7	2.3
-8	2.3
-9	2.0
-10	3.3
-11	1.4



**TRβ-RXR antagonist assay**

LOG[conc.(M)]	luc	ren
-5	-11	4
-6	99	99
-7	110	110
-8	118	110
-9	117	114
-10	113	111

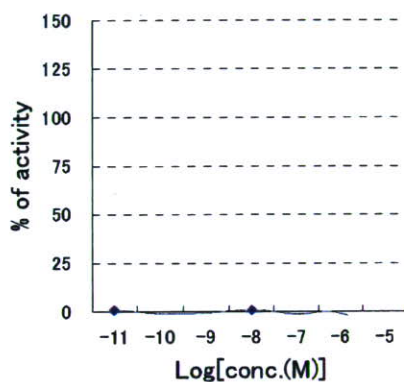




sample No. AT121  
 chemical name 4,4'-THIOBIS(6-TERT-BUTYL-M-CRESOL)  
 CAS. 96-69-5

**AR agonist assay**

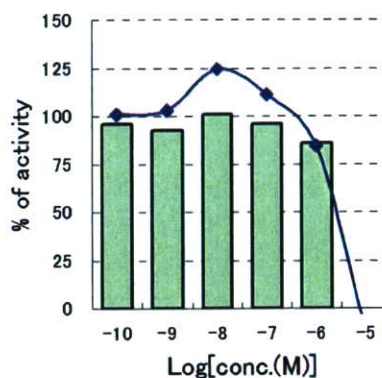
LOG[conc.(M)]	luc
-5	-14.3
-6	-0.8
-7	-1.2
-8	0.9
-9	-0.6
-10	-0.9
-11	0.8



**AR antagonist assay**

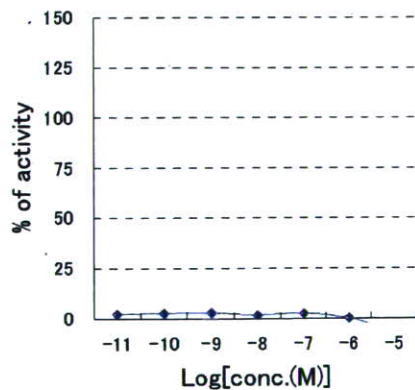
LOG[conc.(M)]	luc	ren
-5	-13	-24
-6	85	86
-7	111	96
-8	124	101
-9	103	92
-10	101	96

Cell viability



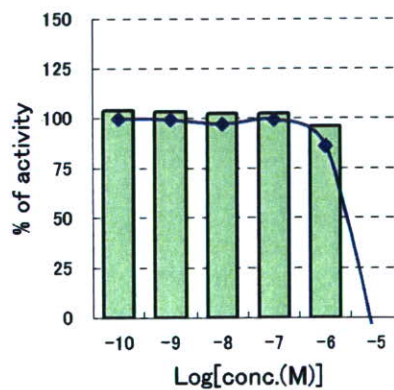
**TRβ-RXR agonist assay**

LOG[conc.(M)]	luc
-5	-6.5
-6	0.7
-7	2.8
-8	2.1
-9	3.1
-10	2.9
-11	2.5



**TRβ-RXR antagonist assay**

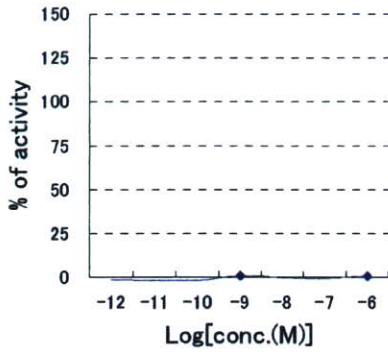
LOG[conc.(M)]	luc	ren
-5	-13	0
-6	86	96
-7	99	102
-8	97	102
-9	99	103
-10	100	104



sample No. AT122  
 chemical name TRIPHENYLPHOSPHINE  
 CAS. 603-35-0

**AR agonist assay**

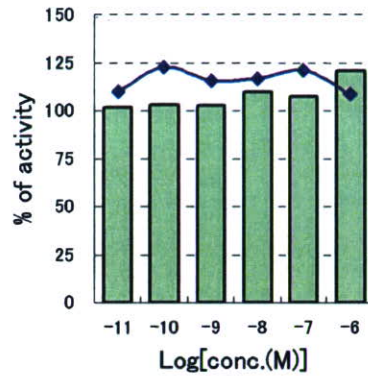
LOG[conc.(M)]	luc
-6	0.5
-7	-0.6
-8	-0.3
-9	1.0
-10	-1.9
-11	-1.6
-12	-1.5



**AR antagonist assay**

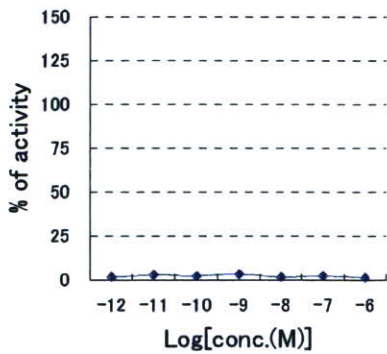
LOG[conc.(M)]	luc	ren
-6	109	121
-7	121	108
-8	117	110
-9	116	103
-10	123	103
-11	110	101

Cell viability



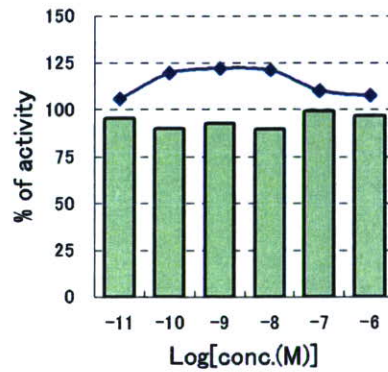
**TRβ-RXR agonist assay**

LOG[conc.(M)]	luc
-6	1.6
-7	2.6
-8	2.0
-9	3.8
-10	2.6
-11	3.2
-12	1.9



**TRβ-RXR antagonist assay**

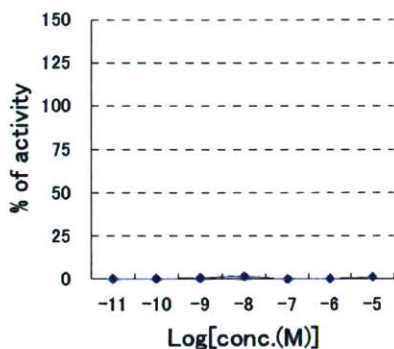
LOG[conc.(M)]	luc	ren
-6	108	97
-7	110	99
-8	121	89
-9	122	93
-10	120	90
-11	106	95



sample No. AT123  
 chemical name 2-BUTANONE OXIME  
 CAS. 96-29-7

**AR agonist assay**

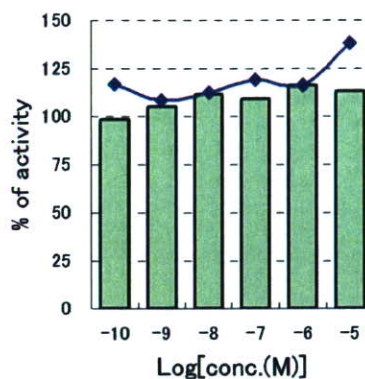
LOG[conc.(M)]	luc
-5	1.4
-6	0.4
-7	0.1
-8	1.8
-9	0.8
-10	0.4
-11	0.2



**AR antagonist assay**

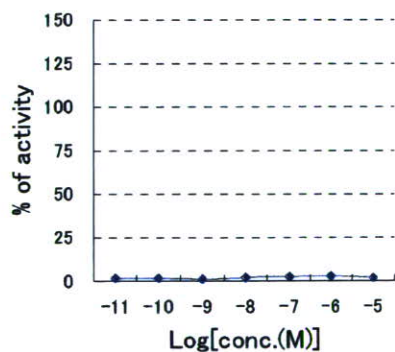
LOG[conc.(M)]	luc	ren
-5	138	113
-6	116	116
-7	119	109
-8	112	111
-9	108	105
-10	117	98

Cell viability



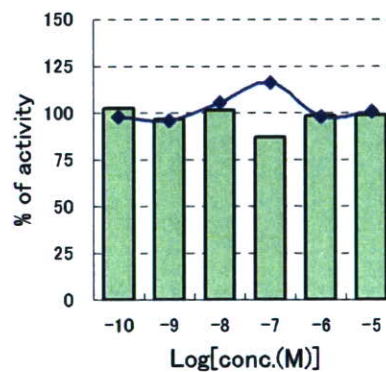
**TRβ-RXR agonist assay**

LOG[conc.(M)]	luc
-5	1.9
-6	3.0
-7	2.8
-8	2.4
-9	1.3
-10	2.1
-11	1.9



**TRβ-RXR antagonist assay**

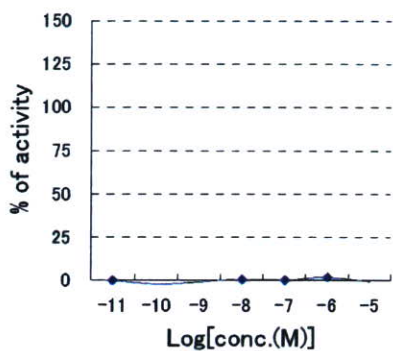
LOG[conc.(M)]	luc	ren
-5	101	99
-6	98	98
-7	116	87
-8	105	102
-9	96	97
-10	98	103



sample No. AT124  
 chemical name 2,2'-Stilbenedisulfonic acid, 4,4'-diamino-  
 CAS. 81-11-8

**AR agonist assay**

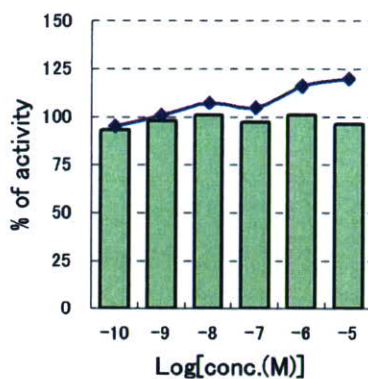
LOG[conc.(M)]	luc
-5	-0.6
-6	2.1
-7	0.5
-8	0.6
-9	-1.0
-10	-2.0
-11	-0.1



**AR antagonist assay**

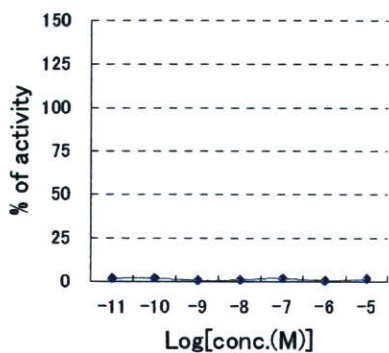
LOG[conc.(M)]	luc	ren
-5	120	96
-6	116	101
-7	105	97
-8	107	101
-9	101	98
-10	95	93

Cell viability



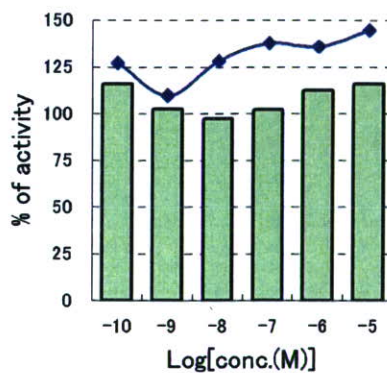
**TRβ-RXR agonist assay**

LOG[conc.(M)]	luc
-5	2.0
-6	1.0
-7	2.2
-8	1.3
-9	1.2
-10	2.4
-11	2.0



**TRβ-RXR antagonist assay**

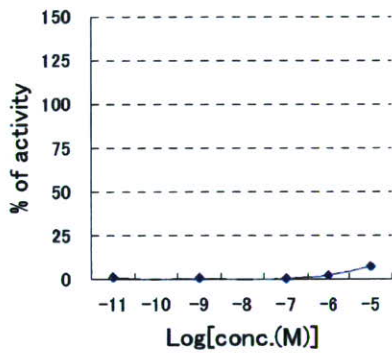
LOG[conc.(M)]	luc	ren
-5	145	116
-6	136	112
-7	138	102
-8	128	97
-9	110	102
-10	127	116



sample No. AT125  
 chemical name Isopropyl 4-hydroxybenzoate  
 CAS. 4191-73-5

**AR agonist assay**

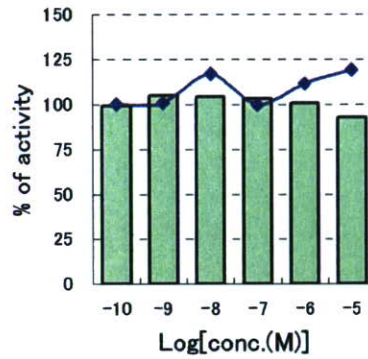
LOG[conc.(M)]	luc
-5	7.6
-6	2.5
-7	0.6
-8	-0.2
-9	1.0
-10	-0.5
-11	1.3



**AR antagonist assay**

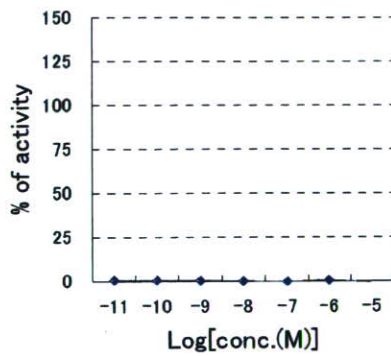
LOG[conc.(M)]	luc	ren
-5	119	93
-6	112	101
-7	100	103
-8	117	104
-9	101	105
-10	100	100

Cell viability



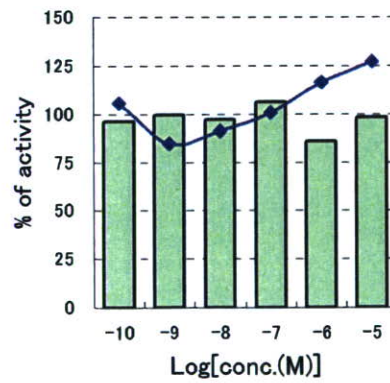
**TRβ-RXR agonist assay**

LOG[conc.(M)]	luc
-5	-0.5
-6	0.7
-7	0.0
-8	0.2
-9	0.6
-10	1.0
-11	1.0



**TRβ-RXR antagonist assay**

LOG[conc.(M)]	luc	ren
-5	127	98
-6	116	86
-7	101	106
-8	91	97
-9	85	100
-10	106	96

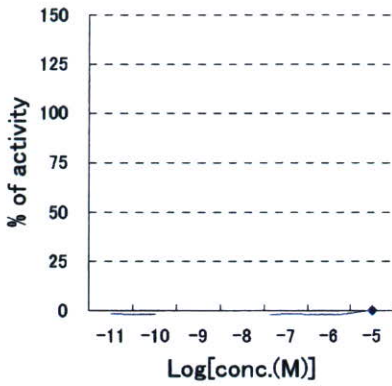




sample No. AT126  
 chemical name Benzidine, 2,2',3,3',5,5',6,6'-octafluoro-  
 CAS. 1038-66-0

**AR agonist assay**

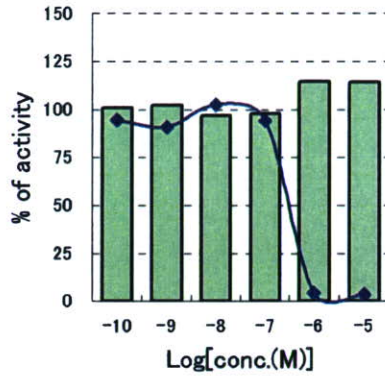
LOG[conc.(M)]	luc
-5	0.2
-6	-2.2
-7	-1.4
-8	-2.8
-9	-1.9
-10	-2.1
-11	-1.4



**AR antagonist assay**

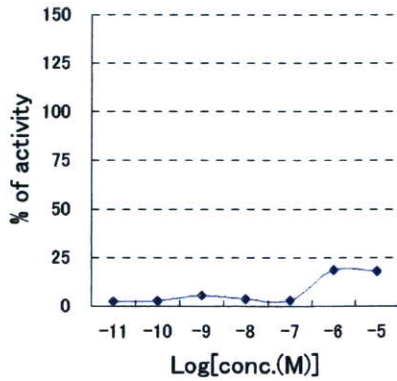
LOG[conc.(M)]	luc	ren
-5	3	114
-6	4	115
-7	94	98
-8	103	97
-9	91	102
-10	94	101

Cell viability



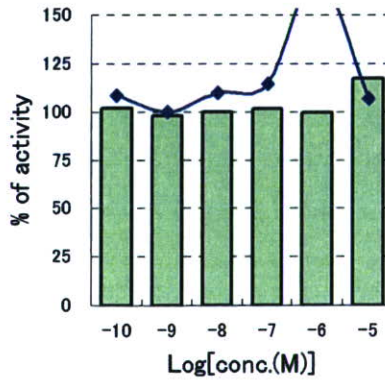
**TRβ-RXR agonist assay**

LOG[conc.(M)]	luc
-5	18.1
-6	18.6
-7	3.2
-8	3.8
-9	5.5
-10	3.0
-11	2.7



**TRβ-RXR antagonist assay**

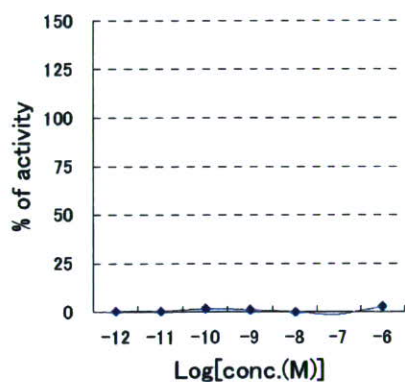
LOG[conc.(M)]	luc	ren
-5	107	117
-6	170	99
-7	115	101
-8	110	100
-9	100	98
-10	109	102



sample No. AT127  
 chemical name Phenol, 4,4'-(diethylideneethylene)di-, diacetate  
 CAS. 84-19-5

**AR agonist assay**

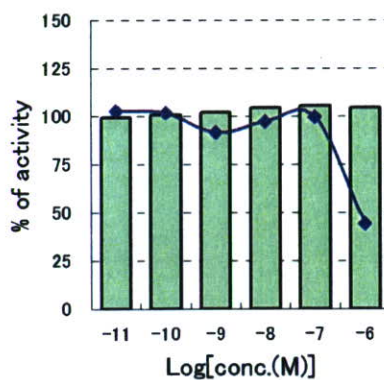
LOG[conc.(M)]	luc
-6	2.9
-7	-1.1
-8	0.0
-9	1.2
-10	1.9
-11	0.3
-12	0.2



**AR antagonist assay**

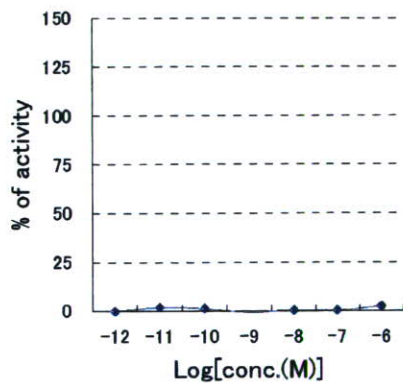
LOG[conc.(M)]	luc	ren
-6	45	105
-7	100	106
-8	97	104
-9	92	102
-10	102	101
-11	103	99

Cell viability



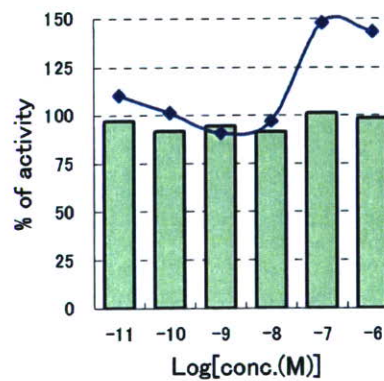
**TRβ-RXR agonist assay**

LOG[conc.(M)]	luc
-6	2.6
-7	0.5
-8	0.6
-9	-0.6
-10	1.6
-11	2.0
-12	0.1



**TRβ-RXR antagonist assay**

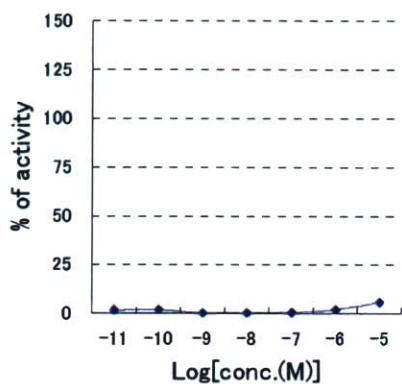
LOG[conc.(M)]	luc	ren
-6	143	98
-7	148	101
-8	97	91
-9	91	94
-10	101	92
-11	110	97



sample No. AT128  
 chemical name Benzidine, 3,3'-dimethyl-  
 CAS. 119-93-7

**AR agonist assay**

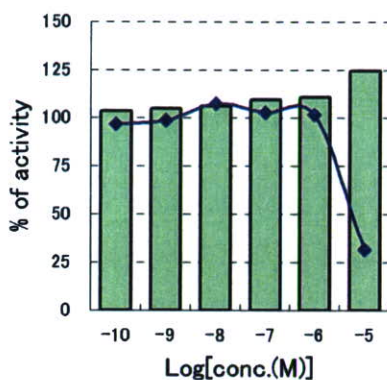
LOG[conc.(M)]	luc
-5	6.0
-6	2.1
-7	0.7
-8	0.3
-9	0.3
-10	2.1
-11	1.7



**AR antagonist assay**

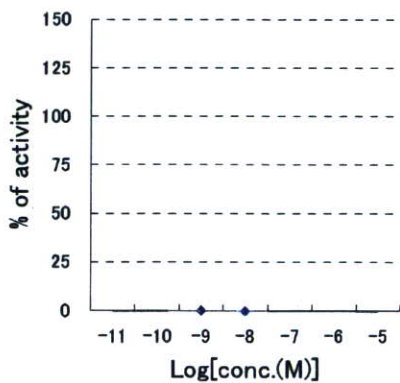
LOG[conc.(M)]	luc	ren
-5	32	124
-6	102	111
-7	103	109
-8	108	106
-9	99	105
-10	97	103

Cell viability



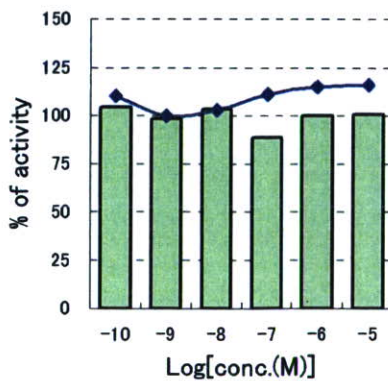
**TRβ-RXR agonist assay**

LOG[conc.(M)]	luc
-5	-0.6
-6	-0.3
-7	-0.3
-8	-0.1
-9	0.4
-10	-0.3
-11	-0.3



**TRβ-RXR antagonist assay**

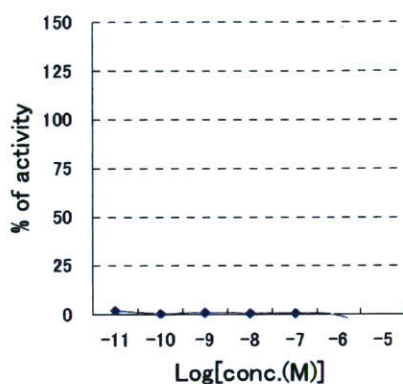
LOG[conc.(M)]	luc	ren
-5	116	101
-6	115	100
-7	111	89
-8	103	103
-9	100	98
-10	110	104



sample No. AT129  
 chemical name 1,2,5,6,9,10-Hexabromocyclododecane  
 CAS. 3194-55-6

**AR agonist assay**

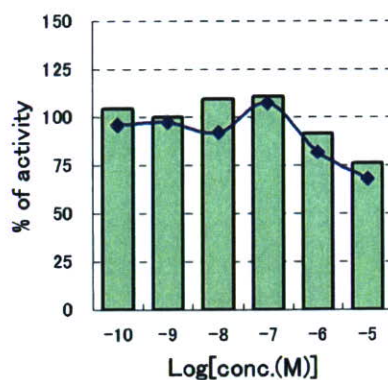
LOG[conc.(M)]	luc
-5	-11.0
-6	-0.5
-7	0.5
-8	0.6
-9	0.9
-10	0.2
-11	2.0



**AR antagonist assay**

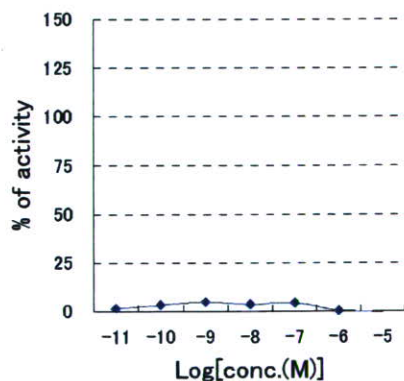
LOG[conc.(M)]	luc	ren
-5	68	76
-6	82	91
-7	108	111
-8	92	110
-9	97	100
-10	96	105

Cell viability



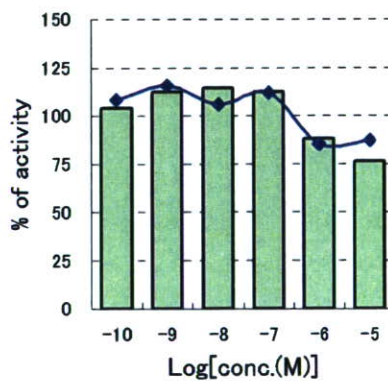
**TRβ-RXR agonist assay**

LOG[conc.(M)]	luc
-5	-0.3
-6	0.5
-7	4.4
-8	4.0
-9	5.1
-10	3.6
-11	1.6



**TRβ-RXR antagonist assay**

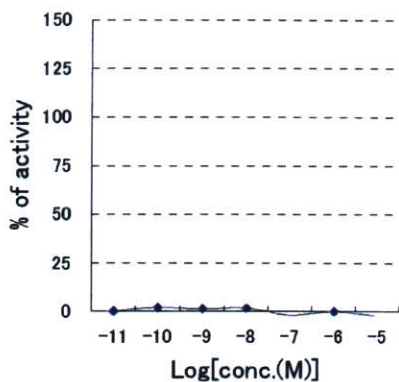
LOG[conc.(M)]	luc	ren
-5	87	76
-6	85	88
-7	112	112
-8	106	114
-9	115	112
-10	109	104



sample No. AT130  
 chemical name Cyclododecane, hexabromo-  
 CAS. 25637-99-4

**AR agonist assay**

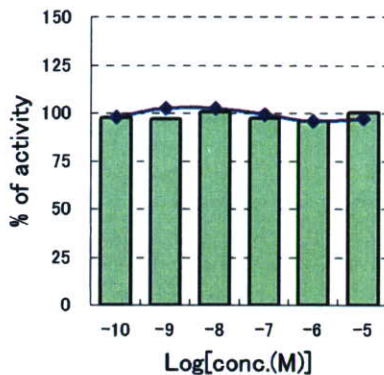
LOG[conc.(M)]	luc
-5	-2.0
-6	-0.1
-7	-1.6
-8	1.9
-9	1.4
-10	2.0
-11	0.2



**AR antagonist assay**

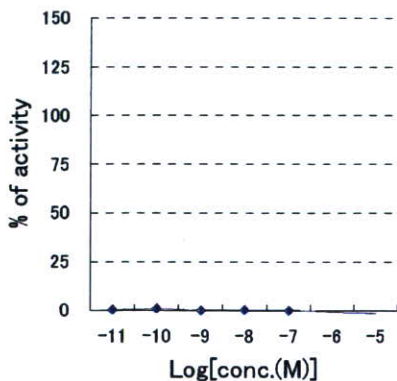
LOG[conc.(M)]	luc	ren
-5	97	100
-6	96	96
-7	100	97
-8	103	101
-9	103	97
-10	98	98

Cell viability



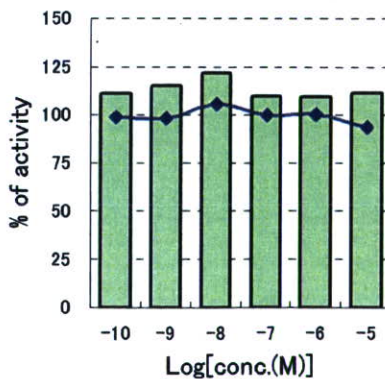
**TRβ-RXR agonist assay**

LOG[conc.(M)]	luc
-5	-1.2
-6	-0.5
-7	0.3
-8	0.5
-9	0.3
-10	1.0
-11	0.6



**TRβ-RXR antagonist assay**

LOG[conc.(M)]	luc	ren
-5	93	111
-6	100	109
-7	100	110
-8	106	122
-9	98	115
-10	99	111

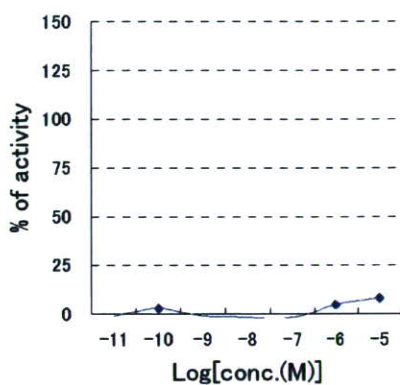




sample No. AT131  
 chemical name Biphenyl, 4,4'-diiodo-  
 CAS. 3001-15-8

**AR agonist assay**

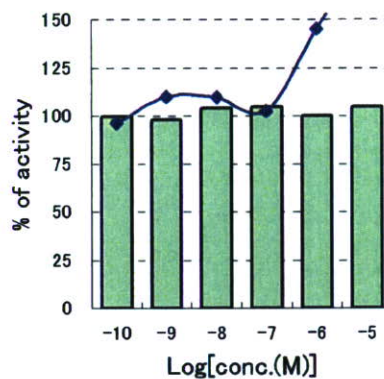
LOG[conc.(M)]	luc
-5	8.2
-6	4.7
-7	-1.6
-8	-1.6
-9	-1.0
-10	2.8
-11	-0.9



**AR antagonist assay**

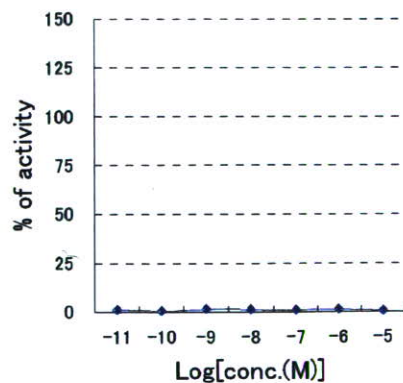
LOG[conc.(M)]	luc	ren
-5	186	104
-6	145	100
-7	102	104
-8	109	104
-9	110	98
-10	96	100

Cell viability



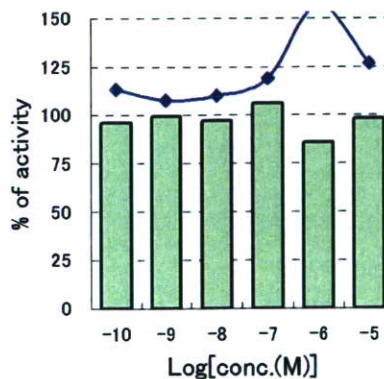
**TRβ-RXR agonist assay**

LOG[conc.(M)]	luc
-5	0.8
-6	1.6
-7	1.2
-8	1.5
-9	1.6
-10	1.0
-11	1.5



**TRβ-RXR antagonist assay**

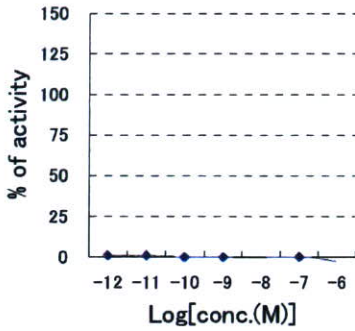
LOG[conc.(M)]	luc	ren
-5	127	98
-6	157	86
-7	119	106
-8	110	97
-9	108	100
-10	113	96



sample No. AT132  
 chemical name 2,2'-Diiodobiphenyl  
 CAS. 2236-52-4

**AR agonist assay**

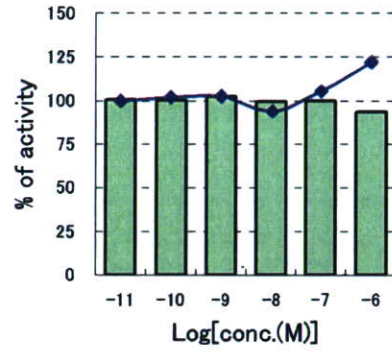
LOG[conc.(M)]	luc
-6	-2.5
-7	0.3
-8	-0.4
-9	0.1
-10	0.2
-11	0.9
-12	1.1



**AR antagonist assay**

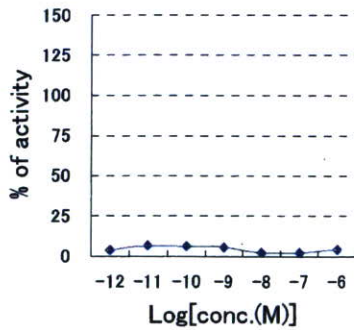
LOG[conc.(M)]	luc	ren
-6	122	94
-7	106	100
-8	94	100
-9	103	102
-10	102	100
-11	100	101

Cell viability



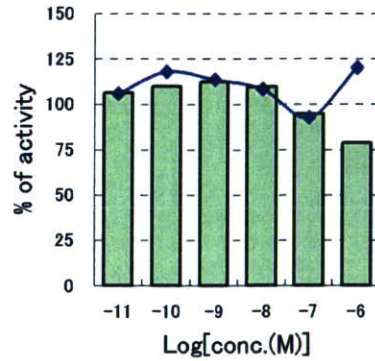
**TRβ-RXR agonist assay**

LOG[conc.(M)]	luc
-6	4.6
-7	2.5
-8	2.5
-9	5.7
-10	6.4
-11	6.9
-12	4.0



**TRβ-RXR antagonist assay**

LOG[conc.(M)]	luc	ren
-6	120	79
-7	93	94
-8	108	110
-9	114	112
-10	118	110
-11	106	106



sample No. AT133